PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 09:42:54 ON 10 MAR 2003 FILE 'CAPLUS' ENTERED AT 09:42:54 ON 10 MAR 2003 COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 1.43 6.14

=> file registry
COST IN U.S. DOLLARS

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 1.84 6.55

FILE 'REGISTRY' ENTERED AT 09:43:39 ON 10 MAR 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 9 MAR 2003 HIGHEST RN 497220-90-3 DICTIONARY FILE UPDATES: 9 MAR 2003 HIGHEST RN 497220-90-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> s dmrie/cn

L13 1 DMRIE/CN

=> d 113

L13 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS

RN 153312-64-2 REGISTRY

CN 1-Propanaminium, N-(2-hydroxyethyl)-N,N-dimethyl-2,3-bis(tetradecyloxy)-,
bromide (9CI) (CA INDEX NAME)

OTHER NAMES:

CN DMRIE

CN N-[1-(2,3-Ditetradecyloxy)propyl]-N,N-dimethyl-N-hydroxyethylammonium bromide

DR 146659-77-0

MF C35 H74 N O3 . Br

CI COM

SR CA

LC STN Files: BIOSIS, CA, CANCERLIT, CAPLUS, IPA, MEDLINE, TOXCENTER, USPATFULL

CRN (191980-81-1)

• Br-

114 REFERENCES IN FILE CA (1962 TO DATE)
6 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
114 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> s dorie/cn L14 2 DORIE/CN

=> d l14 1-2

L14 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2003 ACS

RN 153985-22-9 REGISTRY

CN 1-Propanaminium, N-(2-hydroxyethyl)-N,N-dimethyl-2,3-bis[[(9Z)-1-oxo-9-octadecenyl]oxy]- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 1-Propanaminium, N-(2-hydroxyethyl)-N,N-dimethyl-2,3-bis[(1-oxo-9-octadecenyl)oxy]-, (Z,Z)-

OTHER NAMES:

CN DORIE

FS STEREOSEARCH

DR 139191-30-3

MF C43 H82 N O5

CI COM

SR CA

LC STN Files: CA, CAPLUS, TOXCENTER, USPATFULL

Double bond geometry as shown.

Me Me Me (CH_2) 7 Z (CH_2) 7 Z (CH_2) 7 Z (CH_2) 7 Z (CH_2) 7

PAGE 1-B

__ Me

```
8 REFERENCES IN FILE CA (1962 TO DATE)
               1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
               8 REFERENCES IN FILE CAPLUS (1962 TO DATE)
L14 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2003 ACS
RN
     153312-60-8 REGISTRY
CN
     1-Propanaminium, N-(2-hydroxyethyl)-N, N-dimethyl-2, 3-bis[(9Z)-9-
     octadecenyloxy] -, bromide (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     1-Propanaminium,
N-(2-hydroxyethyl)-N, N-dimethyl-2, 3-bis(9-octadecenyloxy)-
     , bromide, (Z,Z) -
OTHER NAMES:
CN
    DORIE
FS
     STEREOSEARCH
MF
     C43 H86 N O3 . Br
SR
LC
     STN Files: CA, CAPLUS, TOXCENTER, USPATFULL
CRN
     (153985-18-3)
```

Double bond geometry as shown.

Me
$$(CH_2)_7$$
 Z $(CH_2)_8$ O $(CH_2)_8$ Z $(CH_2)_7$ Me Me Me

● Br -

- 4 REFERENCES IN FILE CA (1962 TO DATE)
- 4 REFERENCES IN FILE CAPLUS (1962 TO DATE)

(9CI) (CA INDEX NAME) OTHER CA INDEX NAMES: 1-Propanaminium, N,N,N-trimethyl-2,3-bis(9-octadecenyloxy)-, chloride, (Z,Z)-OTHER NAMES: DOTMA CN CN N-[1-(2,3-Dioleyloxy)propyl]-N,N,N-trimethylammonium chloride STEREOSEARCH 104805-61-0, 375346-70-6 DR C42 H84 N O2 . Cl COM

CI

FS

MF

SR CA

ADISNEWS, AGRICOLA, BIOBUSINESS, BIOSIS, CA, CANCERLIT, LCCAPLUS, CHEMCATS, CIN, MEDLINE, PROMT, TOXCENTER, USPAT2, USPATFULL CRN (122342 - 03 - 4)

Double bond geometry as shown.

Me
$$(CH_2)_7$$
 Z $(CH_2)_8$ O $(CH_2)_8$ Z $(CH_2)_7$ Me $(CH_2)_7$ Me

● cl-

191 REFERENCES IN FILE CA (1962 TO DATE) 14 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 191 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> d l16 1-3

L16 ANSWER 1 OF 3 REGISTRY COPYRIGHT 2003 ACS

RN3700-67-2 REGISTRY

CN 1-Octadecanaminium, N,N-dimethyl-N-octadecyl-, bromide (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Ammonium, dimethyldioctadecyl-, bromide (8CI)

CN Dimethyldioctadecylammonium bromide (6CI, 7CI)

OTHER NAMES:

CN Bis (octadecyl) dimethylammonium bromide

CN DDAB

CNDi-n-octadecyldimethylammonium bromide

CN Dimethyldistearylammonium bromide

CN Dioctadecanyldimethylammonium bromide

CN Dioctadecyldimethylammonium bromide

CN Distearyldimethylammonium bromide

CN DODA (Br)

CN DODAB

CN DSDMAB

GERBU Adjuvant 10 CN

CN GERBU Adjuvant 100

DR 134821-46-8

MF C38 H80 N . Br

CI COM

BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEM, DETHERM*, EMBASE, GMELIN*, IPA, MSDS-OHS, PROMT, TOXCENTER, USPAT2, USPATFULL (*File contains numerically searchable property data) Other Sources: DSL**, EINECS**, TSCA** (**Enter CHEMLIST File for up-to-date regulatory information) CRN (14357-21-2)Me- $(CH_2)_{17} - N^+$ $(CH_2)_{17} - Me$ ● Br-819 REFERENCES IN FILE CA (1962 TO DATE) 25 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 822 REFERENCES IN FILE CAPLUS (1962 TO DATE) 3 REFERENCES IN FILE CAOLD (PRIOR TO 1967) L16 ANSWER 2 OF 3 REGISTRY COPYRIGHT 2003 ACS RN 3282-73-3 REGISTRY CN 1-Dodecanaminium, N-dodecyl-N,N-dimethyl-, bromide (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES: Ammonium, didodecyldimethyl-, bromide (8CI) CNDidodecyldimethylammonium bromide (7CI) OTHER NAMES: CNBis (dodecyl) dimethylammonium bromide CN Bis (n-dodecyl) dimethylammonium bromide CNCNDi-n-dodecyldimethylammonium bromide CN Dilauryldimethylammonium bromide CN Dimethyldidodecylammonium bromide CNDimethyldilaurylammonium bromide MF C26 H56 N . Br CI COM LC STN Files: ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS. BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CEN, CHEMCATS, CHEMLIST, CSCHEM, DETHERM*, GMELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MSDS-OHS, PROMT, TOXCENTER, USPATFULL (*File contains numerically searchable property data) DSL**, EINECS**, TSCA** Other Sources: (**Enter CHEMLIST File for up-to-date regulatory information) CRN (13146-86-6)

ADISNEWS, AGRICOLA, BEILSTEIN*, BIOBUSINESS, BIOSIS,

LC

STN Files:

$$\begin{array}{c} & \text{Me} \\ & | \\ \text{Me}- \text{(CH}_2)_{11} - \text{N}^{+} \text{(CH}_2)_{11} - \text{Me} \\ & | \\ & | \\ & \text{Me} \end{array}$$

● Br-

692 REFERENCES IN FILE CA (1962 TO DATE)

23 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

694 REFERENCES IN FILE CAPLUS (1962 TO DATE)

1 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L16 ANSWER 3 OF 3 REGISTRY COPYRIGHT 2003 ACS

RN 2390-68-3 REGISTRY

CN 1-Decanaminium, N-decyl-N,N-dimethyl-, bromide (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES:

CN Ammonium, didecyldimethyl-, bromide (8CI)

CN Didecyl dimethyl ammonium bromide (6CI, 7CI)

OTHER NAMES:

CN Bromosept

CN DDAB

CN Deciquam

CN Deciquam 222

MF C22 H48 N . Br

CI COM

LC STN Files: ADISNEWS, AGRICOLA, BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CSCHEM, GMELIN*, IFICDB, IFIPAT, IFIUDB, PROMT, TOXCENTER, USPAT2, USPATFULL (*File contains numerically searchable property data)

Other Sources: EINECS**

(**Enter CHEMLIST File for up-to-date regulatory information)
CRN (20256-56-8)

Me
$$(CH_2)_9 - N^+ (CH_2)_9 - Me$$
Me Me

● Br-

214 REFERENCES IN FILE CA (1962 TO DATE)

9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

214 REFERENCES IN FILE CAPLUS (1962 TO DATE)

12 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> s dope/cn

L17 2 DOPE/CN

=> d 117 1-2

```
L17 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2003 ACS
RN
     4004-05-1 REGISTRY
CN
     9-Octadecenoic acid (9Z)-,
(1R) -1-[[[(2-aminoethoxy)hydroxyphosphinyl]oxy]
     methyl]-1,2-ethanediyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     9-Octadecenoic acid (Z)-,
1-[[[(2-aminoethoxy)hydroxyphosphinyl]oxy]methyl
     ]-1,2-ethanediyl ester, (R)-
CN
     Ethanol, 2-amino-, dihydrogen phosphate (ester), monoester with
     1,2-diolein, L- (8CI)
CN
     Olein, 1,2-di-, 2-aminoethyl hydrogen phosphate, L- (8CI)
OTHER NAMES:
CN
     1,2-Dioleoyl-sn-glycero-3-phosphatidylethanolamine
CN
     1,2-Dioleoyl-sn-glycero-3-phosphoethanolamine
CN
     1,2-Dioleoyl-sn-glycero-3-phosphoethanolamine
CN
     1,2-Dioleoyl-sn-glycero-3-phosphorylethanolamine
CN
     Dioleoyl-L-.alpha.-phosphatidylethanolamine
CN
     Dioleoylphosphatidylethanolamine
CN
     DOPE
CN
     L-.alpha.-Dioleoylphosphatidylethanolamine
FS
     STEREOSEARCH
DR
     143062-82-2
MF
     C41 H78 N O8 P
CI
     COM
LC
     STN Files:
                  AGRICOLA, BEILSTEIN*, BIOBUSINESS, BIOSIS, CA, CAOLD,
CAPLUS,
       CASREACT, CHEMCATS, CSCHEM, PROMT, TOXCENTER, USPATFULL
         (*File contains numerically searchable property data)
```

Absolute stereochemistry.
Double bond geometry as shown.

PAGE 1-B

__Me

```
530 REFERENCES IN FILE CAPLUS (1962 TO DATE)
               2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
L17 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2003 ACS
RN
     2462-63-7 REGISTRY
CN
     9-Octadecenoic acid (9Z)-,
1-[[[(2-aminoethoxy)hydroxyphosphinyl]oxy]methy
     l]-1,2-ethanediyl ester (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
     9-Octadecenoic acid (Z)-,
1-[[[(2-aminoethoxy)hydroxyphosphinyl]oxy]methyl
     ]-1,2-ethanediyl ester
CN
     Ethanol, 2-amino-, dihydrogen phosphate (ester), monoester with
     1,2-diolein (8CI)
CN
     Olein, 1,2-di-, 2-aminoethyl hydrogen phosphate (8CI)
     Olein, 1,2-di-, dihydrogen phosphate, 2-aminoethyl ester (7CI)
CN
CN
     Olein, 1,2-di-, phosphate, 2-aminoethyl ester (6CI)
OTHER NAMES:
CN
     1,2-Dioleoyl phosphatidyl ethanolamine
CN
     Dioleoyl (glycerophospho) ethanolamine
CN
     Dioleoyl phosphatidylethanolamine
CN
     DL-Dioleoylphosphatidylethanolamine
CN
     DOPE
CN
     LipofectACE
FS
     STEREOSEARCH
DR
     159317-98-3, 5683-54-5
MF
     C41 H78 N O8 P
CI
     COM
                  AGRICOLA, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA,
LC
     STN Files:
       CANCERLIT, CAOLD, CAPLUS, CSCHEM, EMBASE, IPA, MEDLINE, PROMT,
       TOXCENTER, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
```

524 REFERENCES IN FILE CA (1962 TO DATE)

15 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

Double bond geometry as shown.

PAGE 1-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

```
955 REFERENCES IN FILE CA (1962 TO DATE)
47 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
956 REFERENCES IN FILE CAPLUS (1962 TO DATE)
19 REFERENCES IN FILE CAOLD (PRIOR TO 1967)
```

```
=> s dspe/cn
L18
             0 DSPE/CN
=> s dspe
            13 DSPE
L19
=> s 119 and peg
           196 PEG
L20
             4 L19 AND PEG
=> d 1-4
L20 ANSWER 1 OF 4 REGISTRY COPYRIGHT 2003 ACS
RN
     319491-61-7 REGISTRY
CN
     Poly(oxy-1,2-ethanediyl),
.alpha.-[4-[4-(2,5-dihydro-2,5-dioxo-1H-pyrrol-1-
     yl)phenyl]-1-oxobutyl]-.omega.-[[7-hydroxy-7-oxido-13-oxo-10-[(1-
     oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]oxy]- (9CI)
     (CA INDEX NAME)
OTHER NAMES:
CN
    DSPE-PEG-MPB
MF
     (C2 H4 O)n C57 H97 N2 O12 P
CI
     PMS
PCT Polyether
SR
    CA
LC
     STN Files: CA, CAPLUS
```

$$\begin{array}{c} \text{Me-} & \text{CH}_2)_{16} - \text{C--O} & \text{OH} \\ | & & | \\ \text{Me-} & \text{(CH}_2)_{16} - \text{C--O-CH}_2 - \text{CH--CH}_2 - \text{O--P-O-CH}_2 - \text{CH}_2 - \text{NH--CH}_2 - \text{CH}_2 -$$

PAGE 1-A

$$-O = \begin{bmatrix} CH_2 - CH_2 - O \end{bmatrix}_n C - (CH_2)_3$$

1 REFERENCES IN FILE CA (1962 TO DATE)
1 REFERENCES IN FILE CAPLUS (1962 TO DATE)

L20 ANSWER 2 OF 4 REGISTRY COPYRIGHT 2003 ACS

RN 208044-76-2 REGISTRY

CN Poly(oxy-1,2-ethanediyl), .alpha.-[2-(2,5-dihydro-2,5-dioxo-1H-pyrrol-1-yl)ethyl]-.omega.-[[7-hydroxy-7-oxido-13-oxo-10-[(1-oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]oxy]- (9CI) (CA INDEX NAME) OTHER NAMES:

CN DSPE-PEG-maleimide

DR 441788-07-4, 444083-99-2

MF (C2 H4 O)n C49 H91 N2 O11 P

CI PMS

PCT Polyether

SR CA

LC STN Files: CA, CAPLUS, TOXCENTER, USPATFULL

PAGE 1-A

$$CH_2-CH_2$$
 $O-CH_2-CH_2$ $O-CH_2-CH_2-NH-CH_2-CH_2-O-CH_2-CH_2-NH-CH_2-CH_2-O-CH_2-CH_2-O-CH_2-CH_2-O-CH_$

PAGE 1-B

OH O— C—
$$(CH_2)_{16}$$
— Me $|$ — P— O— CH_2 — CH — CH_2 — O— C — $(CH_2)_{16}$ — Me $|$ 0

12 REFERENCES IN FILE CA (1962 TO DATE)

3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 12 REFERENCES IN FILE CAPLUS (1962 TO DATE)

L20 ANSWER 3 OF 4 REGISTRY COPYRIGHT 2003 ACS RN 170931-04-1 REGISTRY

CN Poly(oxy-1,2-ethanediyl), .alpha.-[(10R)-7-hydroxy-7-oxido-13-oxo-10-[(1-oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]-.omega.-hydroxy- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Poly(oxy-1,2-ethanediy1), .alpha.-[7-hydroxy-7-oxido-13-oxo-10-[(1-oxooctadecy1)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-y1]-.omega.-hydroxy-, (R)-

OTHER NAMES

CN 1,2-Distearoyl-sn-glycero-3-phosphoethanolamine-N-[poly(ethylene glycol)]

CN DSPE-PEG

CN PEG-DSPE

CN Poly(oxy-1,2-ethanediyl), .alpha.-[7-hydroxy-13-oxo-10-[(1-oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]-.omega.-hydroxy-, P-oxide, (R)-

MF (C2 H4 O)n C43 H86 N O9 P

CI PMS

PCT Polyether

SR CA

LC STN Files: ADISNEWS, BIOSIS, CA, CAPLUS, TOXCENTER, USPATFULL

PAGE 1-A $Me - (CH_2)_{16} - C - O - CH_2$ OH $CH_2 - CH_2 - O - CH_2 - CH_2 - O - P - O - CH_2 - CH - O - CH_2$

PAGE 1-B

49 REFERENCES IN FILE CA (1962 TO DATE)

1 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

50 REFERENCES IN FILE CAPLUS (1962 TO DATE)

L20 ANSWER 4 OF 4 REGISTRY COPYRIGHT 2003 ACS

RN 145035-96-7 REGISTRY

CN Poly(oxy-1,2-ethanediyl), .alpha.-[7-hydroxy-7-oxido-13-oxo-10-[(1-oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]-.omega.-hydroxy- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Poly(oxy-1,2-ethanediyl), .alpha.-[7-hydroxy-13-oxo-10-[(1-

oxooctadecyl)oxy]-6,8,12-trioxa-3-aza-7-phosphatriacont-1-yl]-.omega.-hydroxy-, P-oxide

OTHER NAMES:

CN Distearoylphosphatidylethanolamine-PEG

CN DSPE-PEG

CN PEG-DSPE

MF (C2 H4 O)n C43 H86 N O9 P

CI PMS, COM

PCT Polyether

SR CA

LC STN Files: BIOSIS, CA, CAPLUS, TOXCENTER, USPAT2, USPATFULL

PAGE 1-2

PAGE 1-B

128 REFERENCES IN FILE CA (1962 TO DATE)

14 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

131 REFERENCES IN FILE CAPLUS (1962 TO DATE)

=> log hold

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY

SESSION

60.40

66.95

SESSION WILL BE HELD FOR 60 MINUTES

L12 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2003 ACS

AN 1996:724086 CAPLUS

DN 125:338691

TI Skin irritation-free stable cosmetics containing amido-containing anionic surfactants

IN Nishio, Hiroyuki; Tsubone, Kazuyuki

PA Kanebo Ltd, Japan

SO Jpn. Kokai Tokkyo Koho, 6 pp.

CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO. DATE

PI JP 08245365 A2 19960924 JP 1995-83305 19950314

PRAI JP 1995-83305 19950314

OS MARPAT 125:338691